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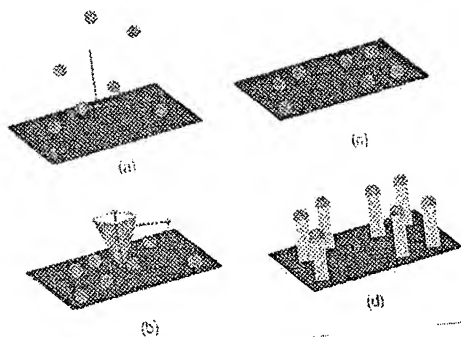


Figure 1

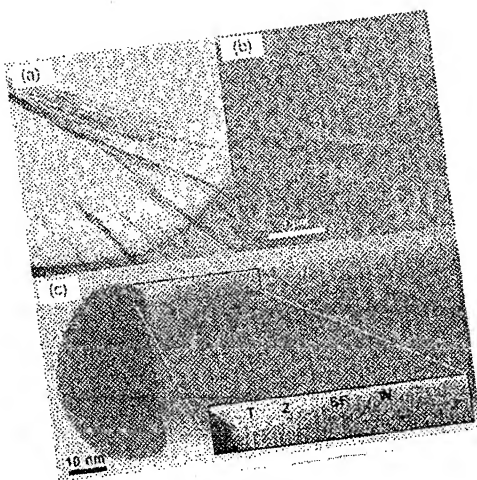
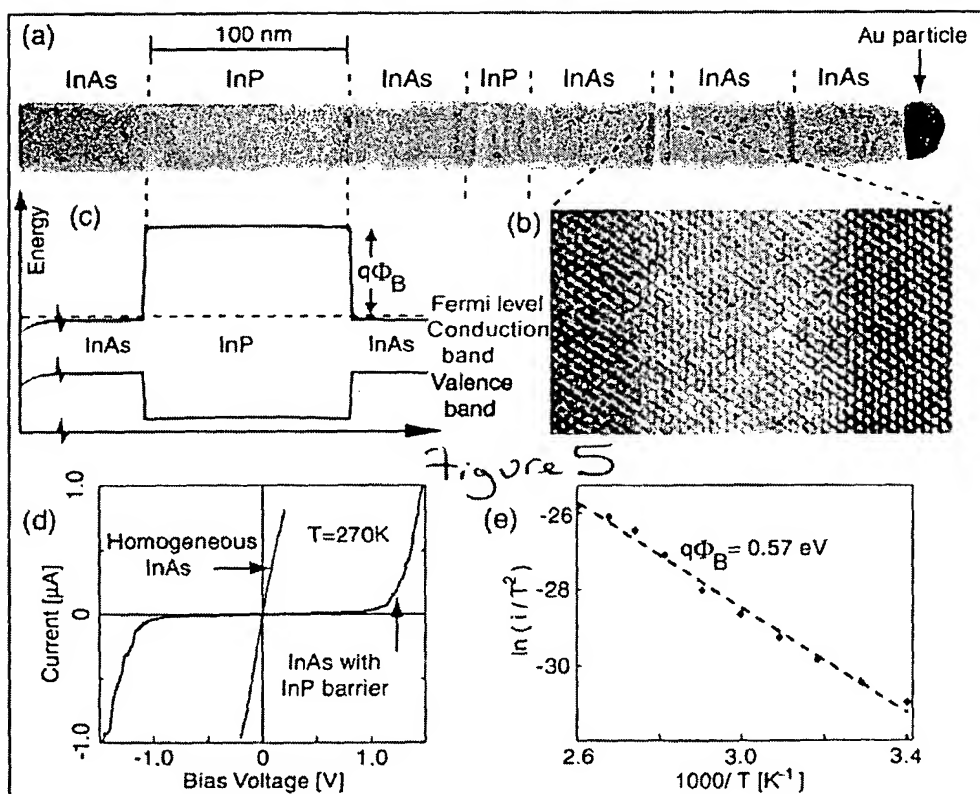
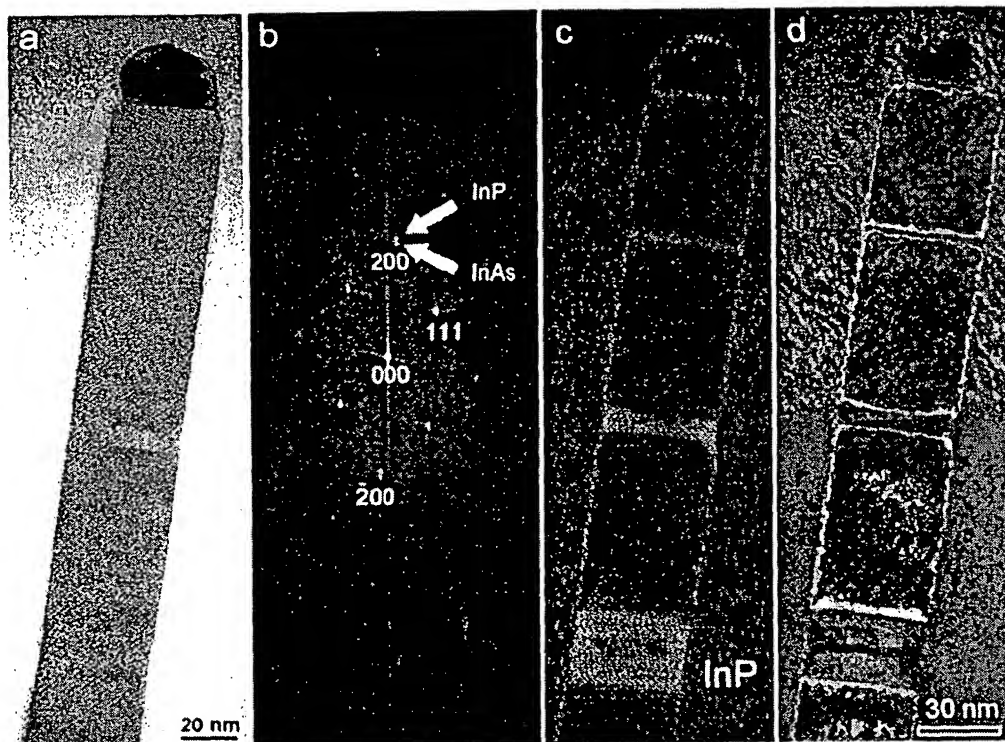


Figure 2

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Figure 4



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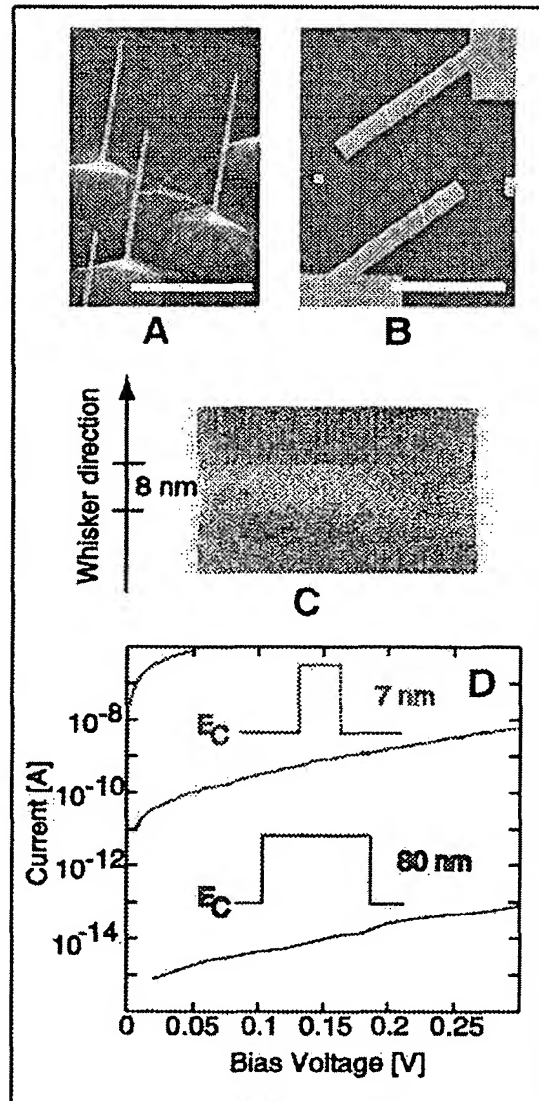


Figure 6

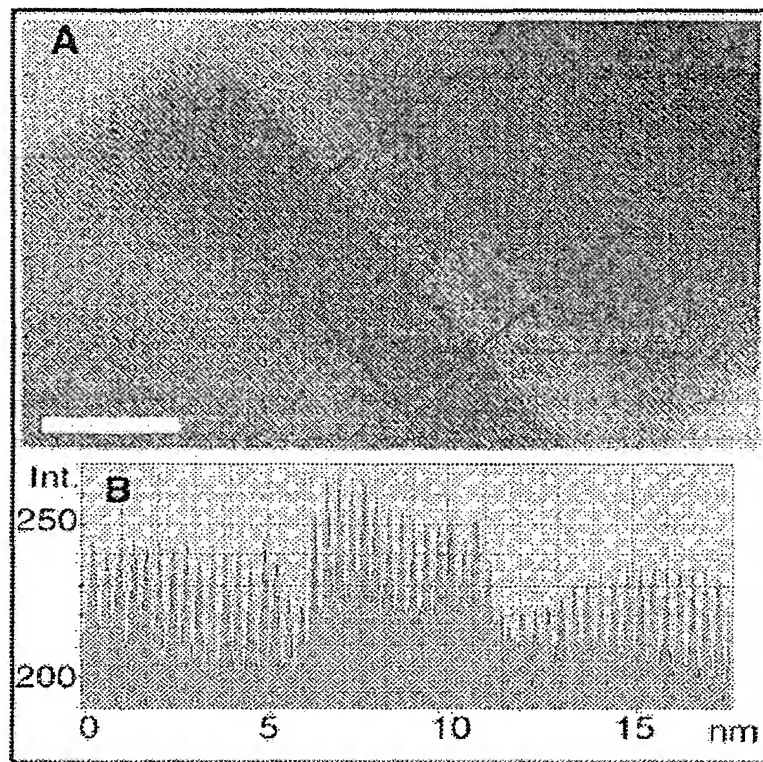


Figure 7

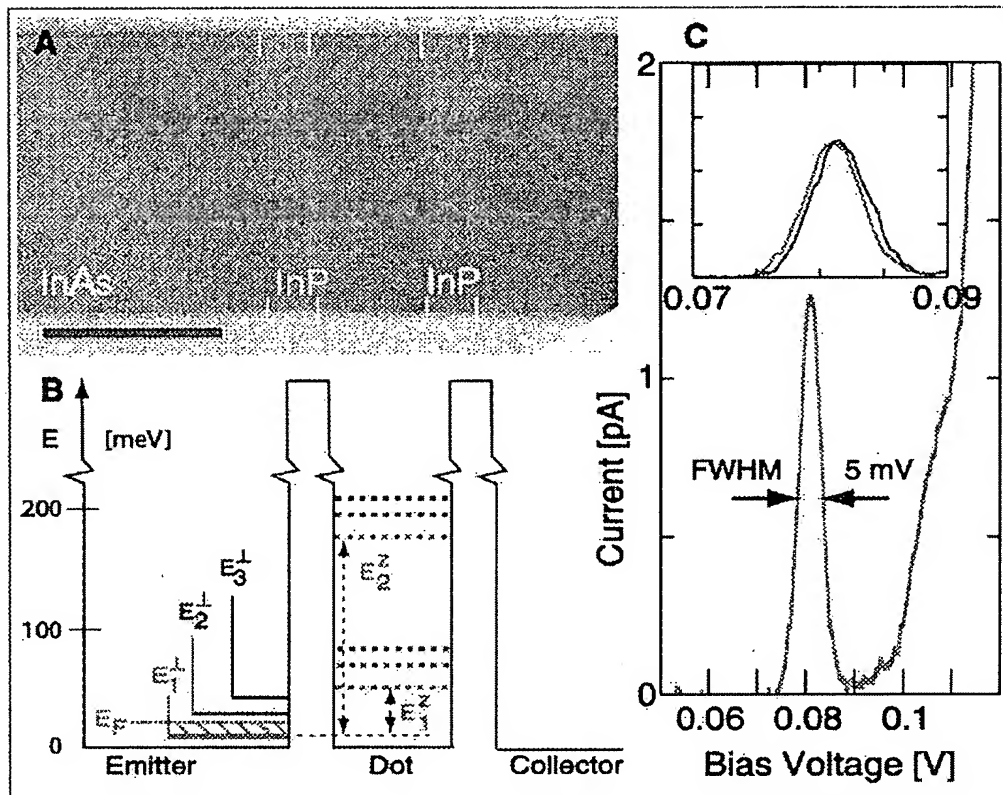
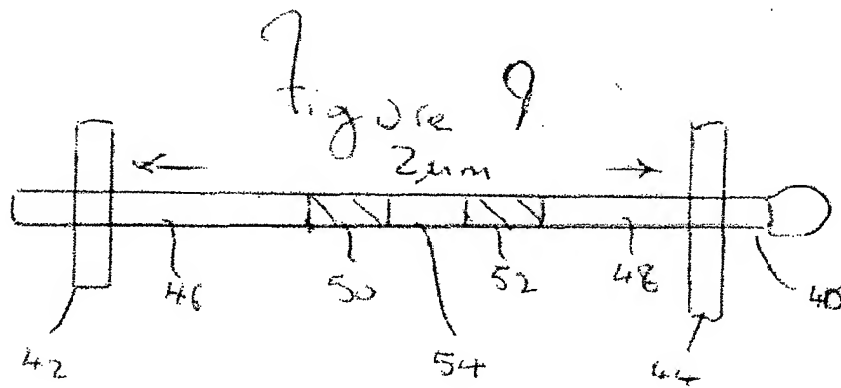
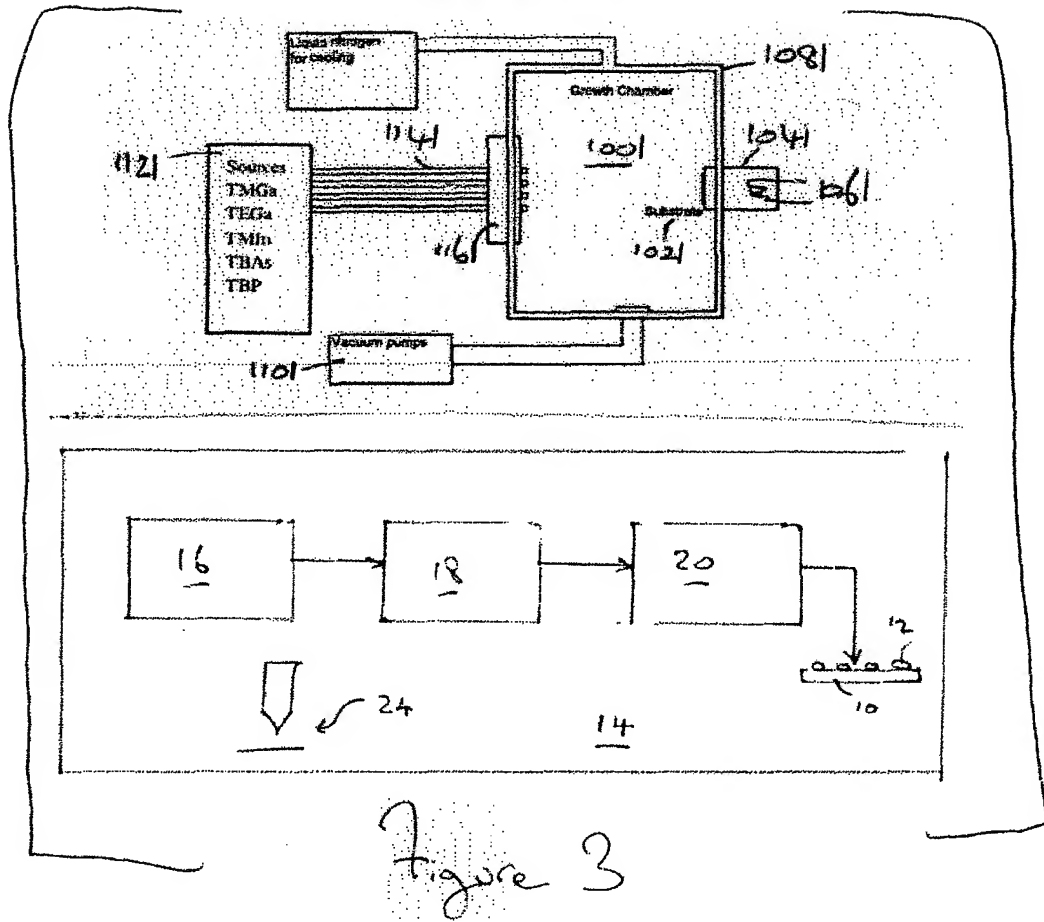


Figure 8

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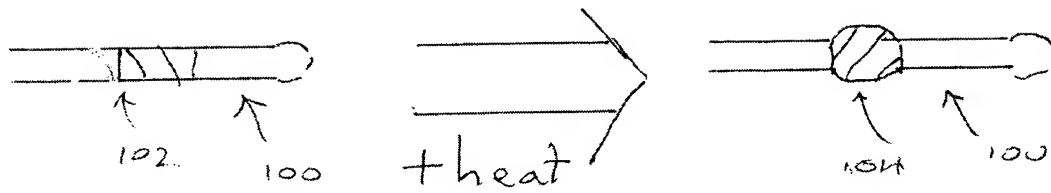


Figure 10

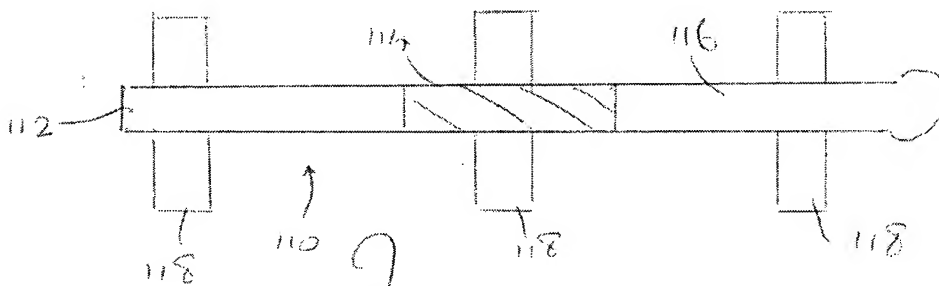
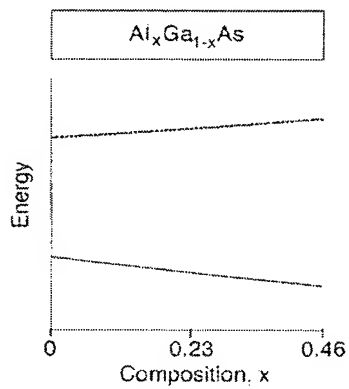


Figure 11



graded semiconductor

Figure 13

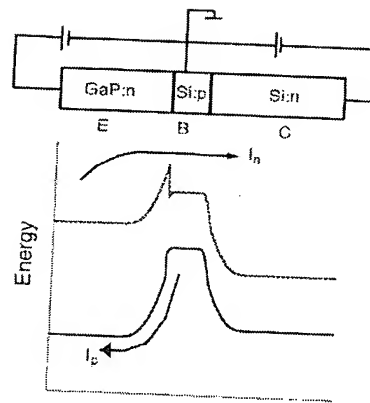


Figure 12

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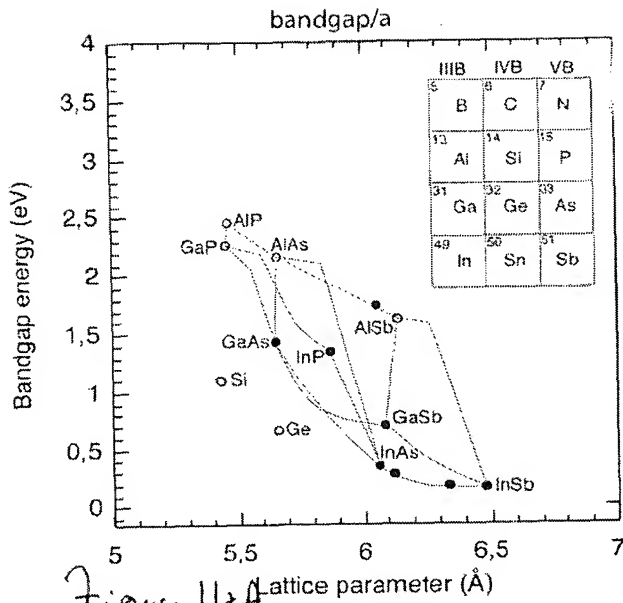


Figure 14A

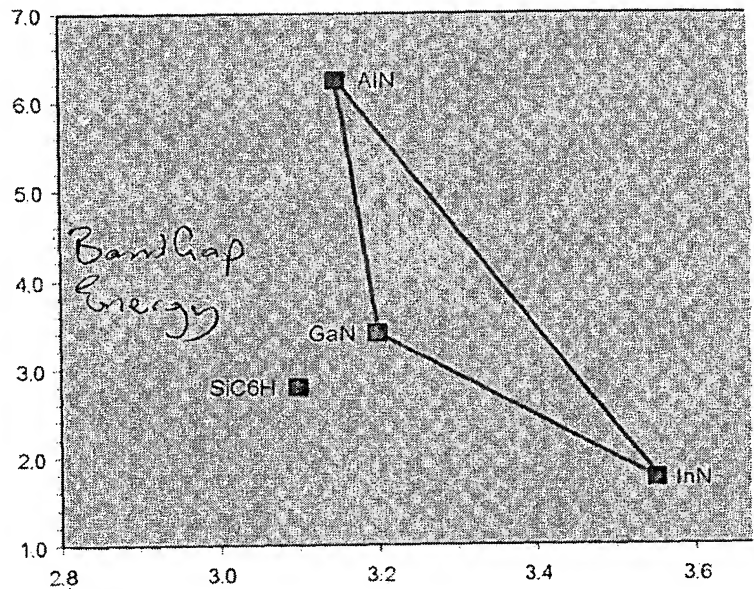


Figure 14B

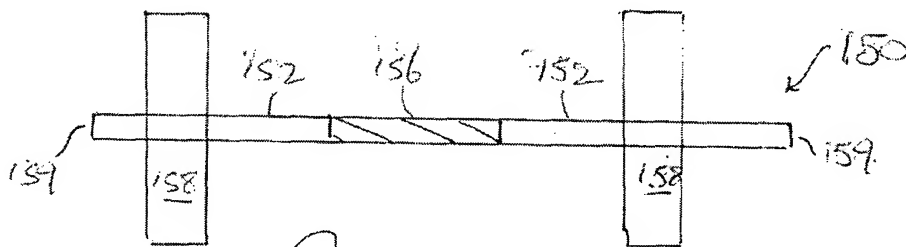


Figure 15

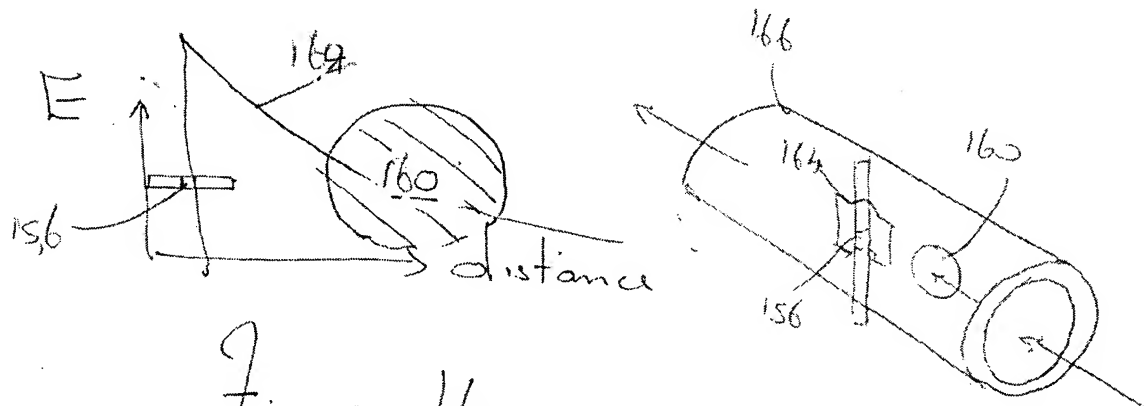


Figure 16

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Figure 17

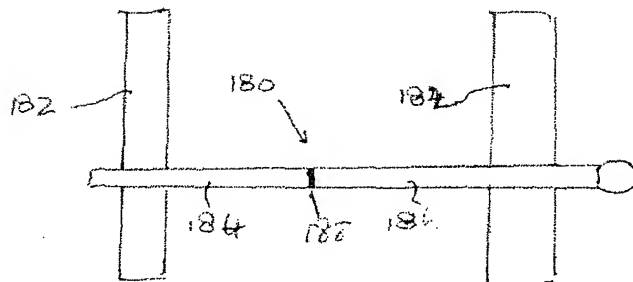
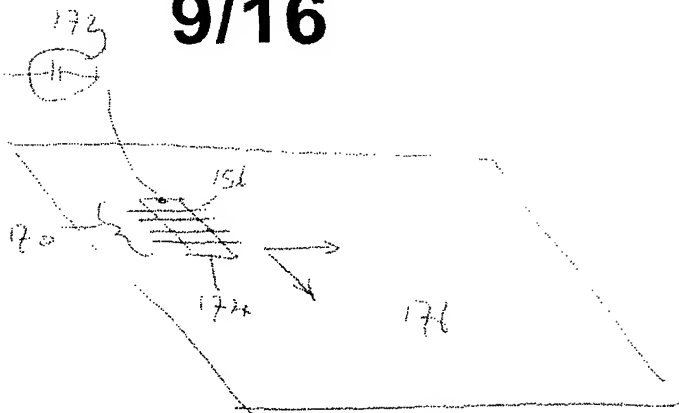
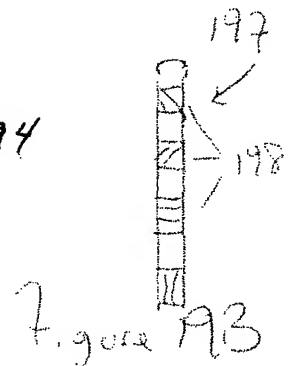
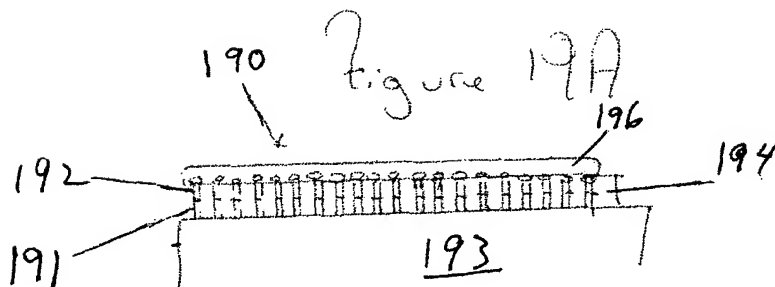
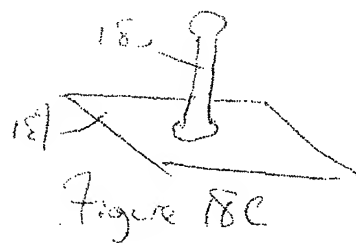
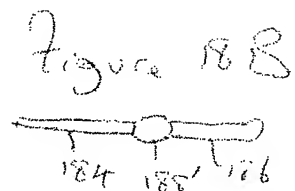


Figure 18A



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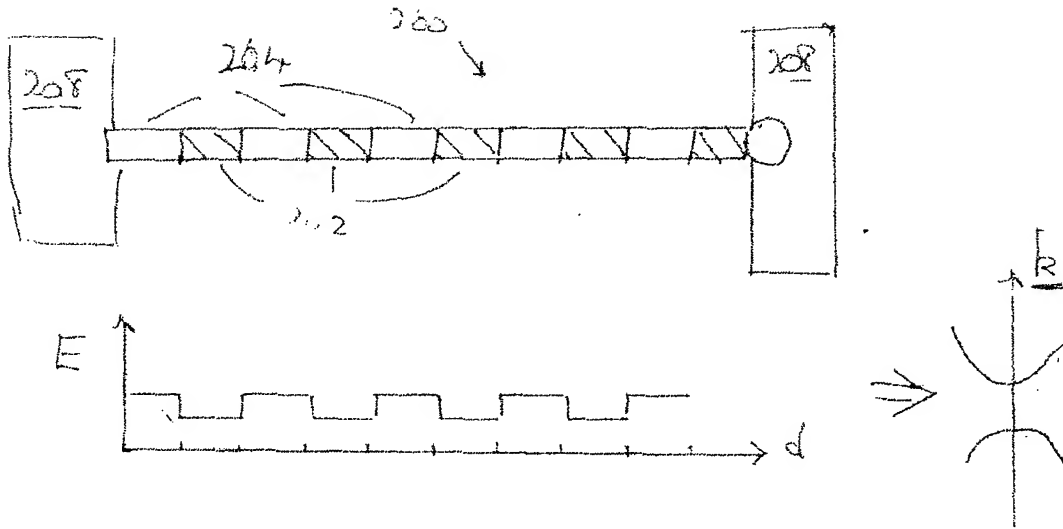


Figure 20

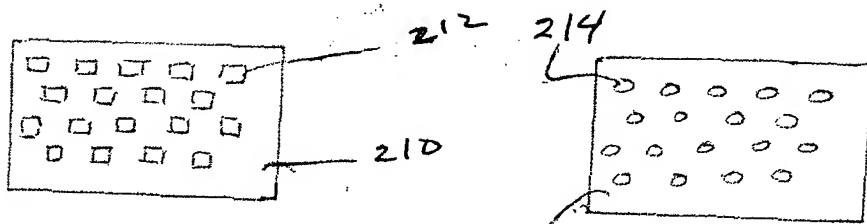


Figure 21A

Figure 21B

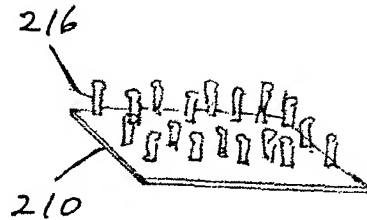
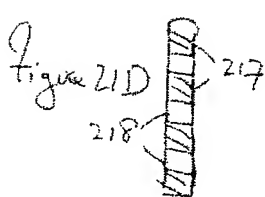


Figure 21C

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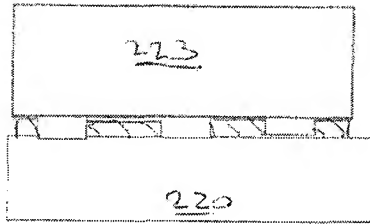


Figure 22A

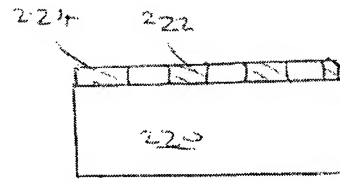


Figure 22B

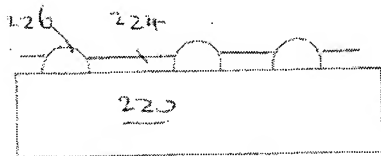


Figure 22C

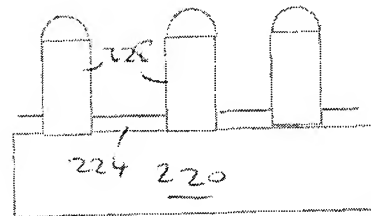


Figure 22D

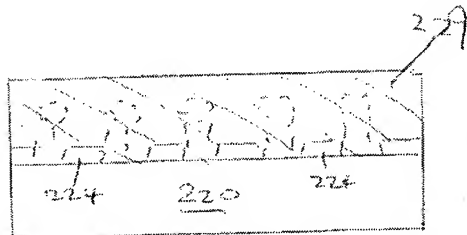


Figure 22E

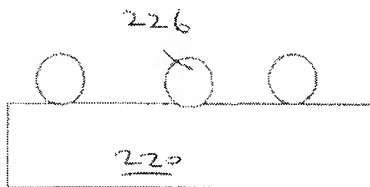


Figure 22F

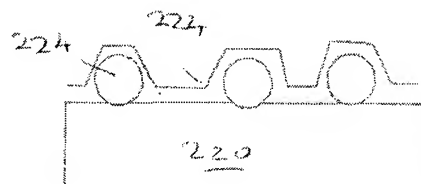


Figure 22G

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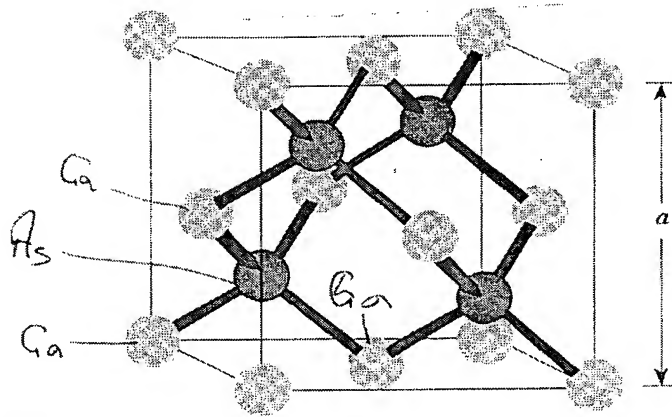
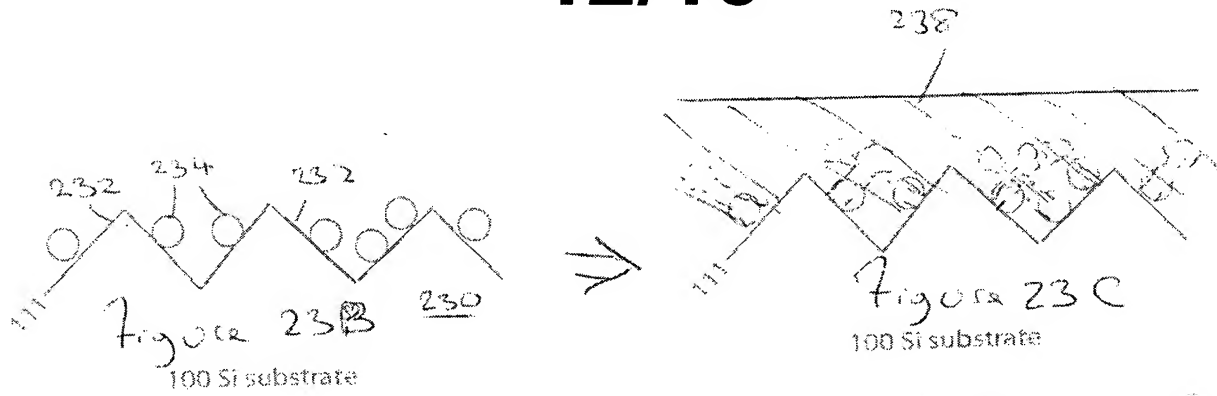
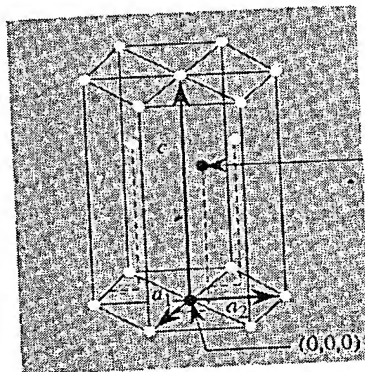
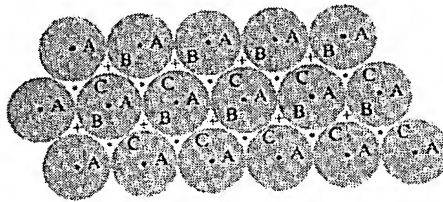


Figure 23A The zinc blende crystal structure. The structure consists of the interpenetrating fcc lattices, one displaced from the other by a distance $(\frac{a}{4}, \frac{a}{4}, \frac{a}{4})$ along the body diagonal; underlying Bravais lattice is fcc with a two atom basis. The positions of the two atoms are $(0,0,0)$ and $(\frac{a}{4}, \frac{a}{4}, \frac{a}{4})$.



$|c| = |a_3| = 1.633 |a_1|$
Two atoms per basis

$(\frac{2}{3}a_1 + \frac{1}{3}a_2 + \frac{1}{2}a_3)$

Figure 24A

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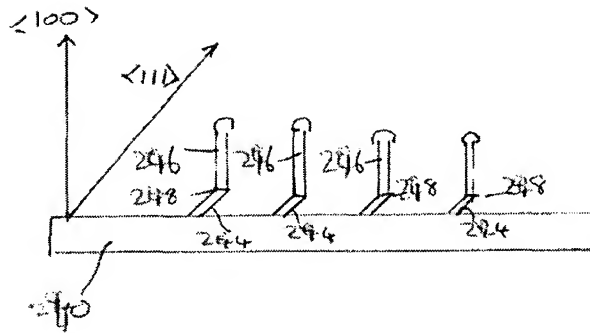


Figure 24B

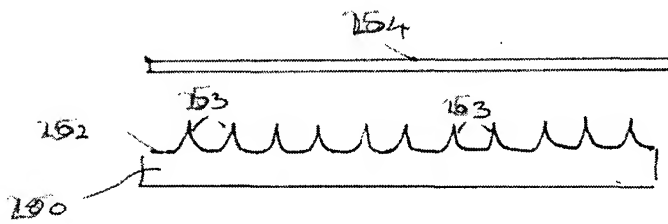


Figure 25A

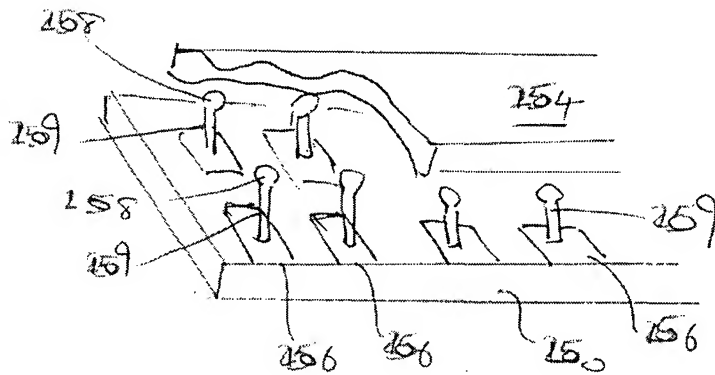


Figure 25B

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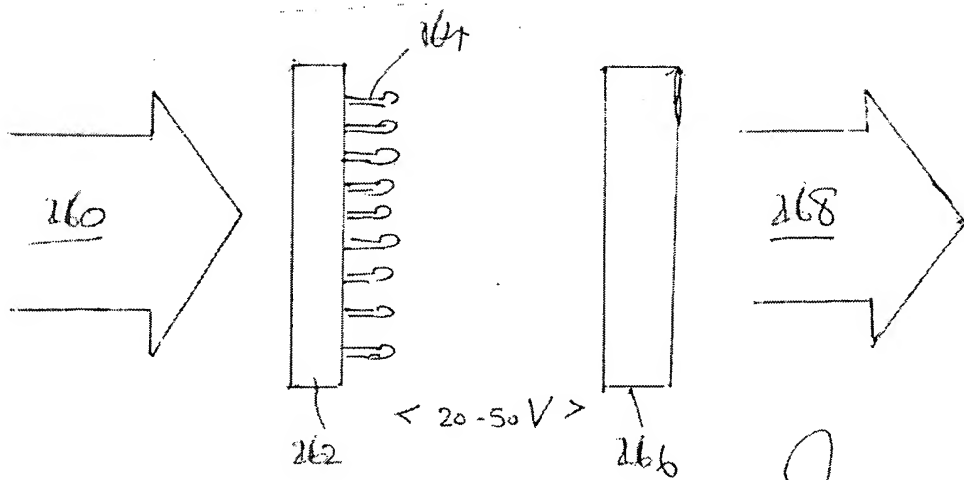


Figure 26

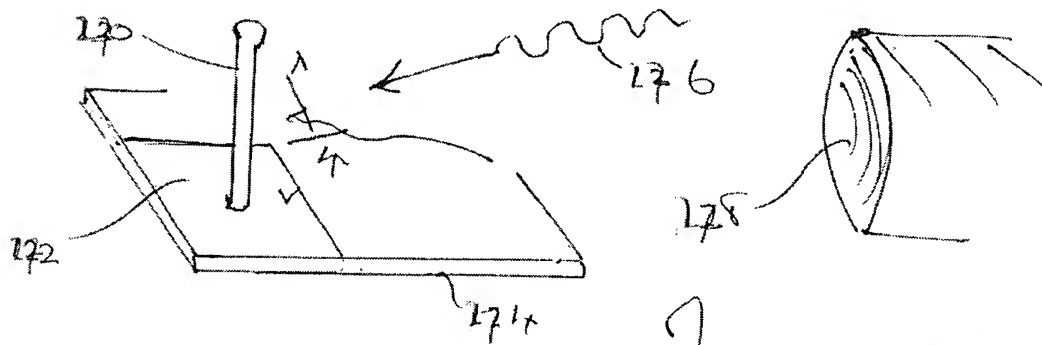


Figure 27

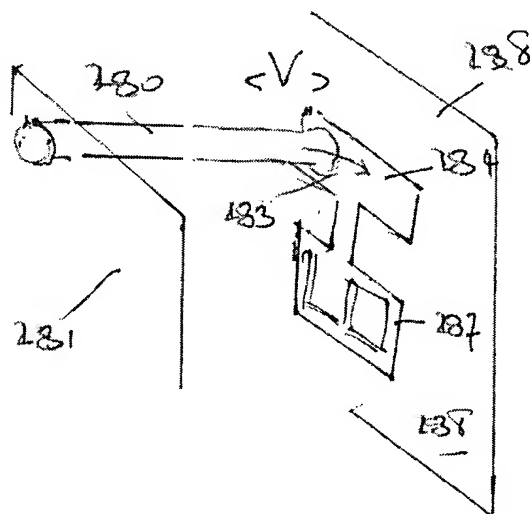


Figure 28

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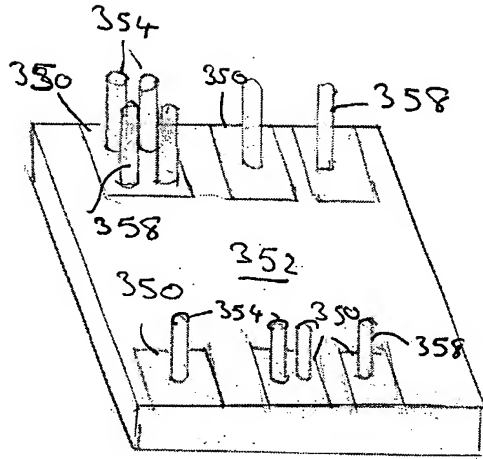


Figure 29

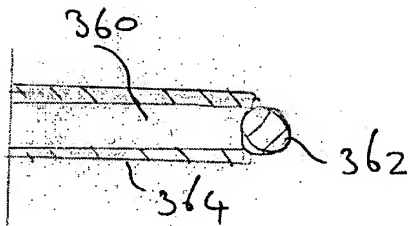


Figure 30

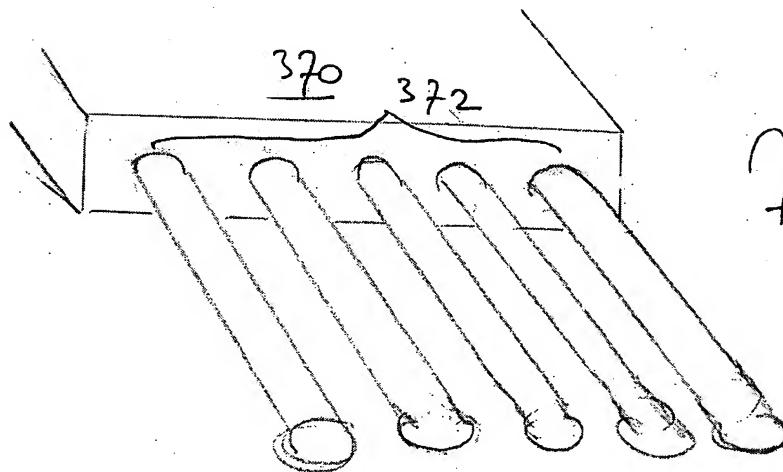


Figure 31

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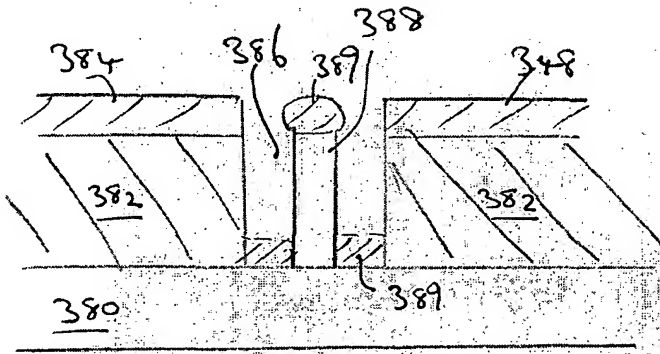


Figure 32

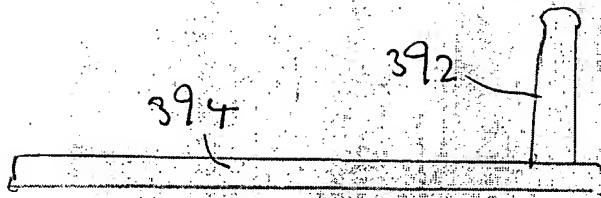


Figure 33